Patent Application Attorney Docket No. PC10973B

IN THE CLAIMS

(Currently amended) A compound of general formula (I), or 1. pharmaceutically acceptable salts, solvates or polymorphs thereof;

wherein:

L and U, which may be the same or different, are -N-, -N*(-O')- or -C(H)-; M and Q, which may be the same or different, are -N-, -N $^+$ (-O $^-$)- or -C(R 4)-; wherein ring A contains 1 nitrogen atom;

R¹ and R², which may be the same or different, are hydrogen, C₁₋₆alkyl, $(CH_2)_m(C_{3^{-8}}$ cycloalkyl) wherein m = 0, 1, 2 or 3, or R¹ and R² together with the nitrogen to which they are attached form an azetidine ring;

W. Y and Z. which may be the same or different, are hydrogen, halogen, C14alkyl, CF3, OCF3, C14alkylthio or C14alkoxy; or Y and Z are linked so that, together with the interconnecting atoms, Y and Z form a fused 5 to 7-membered carbocyclic or heterocyclic ring which may be saturated, unsaturated or aromatic, and wherein when Y and Z form a heterocyclic ring, in addition to carbon atoms, the linkage contains one or two heteroatoms independently selected from oxygen, sulfur and nitrogen; and wherein W, Y and Z are not all hydrogen; and

each R4 is independently:

A-X, wherein A = -(CH₂)₀- where p is 0, 1 or 2; X is hydrogen, CONR⁶R⁷. SO₂NR⁶R⁷, SO₂NHC(=O)R⁶, hydroxy, C₁₋₄alkoxy, NR⁶SO₂R⁹, NO₂, NR⁶R¹¹, CN, CO₂R¹⁰. SR¹⁰, S(O)R⁹ or SO₂R¹⁰; R⁶, R⁷, R⁸ and R¹⁰ which may be the same or different, are hydrogen or C₁₋₆alkyl optionally substituted independently by one or more R¹²; R⁹ is C₁₋₆ alkyl optionally substituted independently by one or more R¹²; R¹¹ is hydrogen, C₁₋₆ alkyl

Patent Application Attorney Docket No. PC10973B

optionally substituted independently by one or more R¹², C(O)R⁶, CO₂R⁹, C(O)NHR⁶ or SO₂NR⁶R⁷; R¹² is fluoro, hydroxy, CO₂H, C_{3.6}cycloalkyl, NH₂, CONH₂, C_{1.6}alkoxy, C₁. 6alkoxycarbonyl or a 5- or 6-membered heterocyclic ring containing 1, 2 or 3 heteroatoms selected from N, S and O optionally substituted independently by one or more R¹³; or R⁶ and R⁷, together with the nitrogen to which they are attached, form a 4-, 5- or 6-membered heterocyclic ring optionally substituted independently by one or more R¹³; or a 5- or 6-membered heterocyclic ring containing 1, 2 or 3 heteroatoms selected from N, S and O, optionally substituted independently by one or more R13; wherein R13 is hydroxy, C₁-C₄alkoxy, fluoro, C₁-C₆alkyl, haloalkyl, haloalkoxy, -NH₂, -NH(C₁-C₆alkyl) or -N(C₁-C₆alkyl)₂.

- 2. (Cancelled)
- (Previously presented) A compound according to claim 1 wherein L is -3. C(H)-.
- (Original) A compound according to claim 1 wherein R¹ and R², which may be the same or different, are hydrogen or C₁-C₆alkyl, or R¹ and R², together with the nitrogen to which they are attached, form an azetidine ring.
- (Original) A compound according claim 1 wherein R¹ is methyl and R² is 5. hydrogen or methyl, or R1 and R2, together with the nitrogen to which they are attached. form an azetidine ring.
- (Original) A compound according to claim 1 wherein R1 is methyl and R2 6. is hydrogen or methyl.
- 7. (Original) A compound according to claim 1 wherein W is hydrogen, C₁₋₈alkyl, C₁₋₄alkoxy or halogen.
- 8. (Original) A compound according to claim 1 wherein W is hydrogen, methyl or ethyl; and Y and Z, which may be the same or different, are hydrogen, methyl, ethyl, CF₃, OCF₃, methylthio, ethylthio, methoxy, ethoxy, chloro, fluoro or bromo; or Y

Patent Application Attorney Docket No. PC10973B

and Z are linked so that, together with the interconnecting atoms, Y and Z form a fused 5 to 7-membered carbocyclic or heterocyclic ring which may be saturated, unsaturated or aromatic, and wherein when Y and Z form a heterocyclic ring, in addition to carbon atoms, the linkage contains one or two heteroatoms independently selected from oxygen, sulfur and nitrogen; wherein W, Y and Z are not all hydrogen.

- 9. (Original) A compound according to claim 1 wherein W is hydrogen; and Y and Z, which may be the same or different, are hydrogen, fluoro, chloro, methyl, ethyl, methylthio, ethylthio, methoxy or ethoxy; or Y and Z are linked so that, together with the interconnecting atoms, Z and Y form a fused 5 to 7-membered heterocyclic ring containing one or more sulfur atoms; wherein Y and Z are not both hydrogen.
- 10. (Original) A compound according to claim 1 wherein when Y and Z are linked so that, together with the interconnecting atoms, Z and Y form a fused 5 to 7-membered heterocyclic ring containing one or more sulfur atoms, the linkages forming the fused ring are -S(CH₂)₂-, -CH₂S-CH₂- or -S(CH₂)₂O- wherein either end of these linkages correspond to either group Y or Z.
- 11. (Original) A compound according to claim 1 wherein, when present, each R^4 is independently -(CH_2)_p-X, where p is 0, 1 or 2; X is hydrogen, $CONR^6R^7$, $SO_2NR^6R^7$, $SO_2NH(C=O)R^6$, hydroxy, C_{1-4} alkoxy, $NR^8SO_2R^9$, NO_2 , NR^6R^{11} , CN, CO_2R^{10} , SR^{10} , $S(O)R^9$ or SO_2R^{10} ; wherein R^8 , R^7 , R^8 , R^{10} or R^{11} , which may be the same or different, are hydrogen or C_{1-6} alkyl; and R^9 is C_{1-6} alkyl.
- 12. (Original) A compound according to claim 1 wherein, when present each R⁴ is independently -(CH₂)_p-X, where p is 0 or 1; X is hydrogen, CONR⁶R⁷, SO₂NR⁶R⁷, NR⁸SO₂R⁹, hydroxy or NR⁶R¹¹; wherein R⁶, R⁷, R⁸, or R¹¹, which may be the same or different, are hydrogen or C₁₋₆alkyl; and R⁹ is C₁₋₆alkyl.
- _13. (Original) A compound according to claim 1 wherein the compound is selected from:

N-methyl-N-({4-[4-(methylsulfanyl)phenoxy]-3-pyridinyl}methyl)amine,
N-{[4-(2,3-dihydro-1-benzothien-5-yloxy)-3-pyridinyl]methyl}-N-methylamine,
N-({4-[3-chloro-4-(methylsulfanyl)phenoxy]-3-pyridinyl}methyl)-N-methylamine,
N-methyl-N-({3-[4-(methylsulfanyl)phenoxy]-4-pyridinyl}methyl)amine,
N-methyl-N-({3-[3-methyl-4-(methylsulfanyl)phenoxy]-4-pyridinyl}-methyl)amine,

and

Patent Application Attorney Docket No. PC10973B

N-{[4-(2,3-Dihydro-1,4-benzoxathiin-7-yloxy)-6-methyl-3-pyridinyl]methyl}-*N*-methylamine,

N-methyl-N-({6-methyl-4-[3-methyl-4-(methylsulfanyl)phenoxy]-3-pyridinyl}methyl)amine,

N-({4-[3-chloro-4-(methylsulfanyl)phenoxy]-3-pyridinyl}methyl)-*N*,*N*-dimethylamine,

N-({4-[3-fluoro-4-(methylsulfanyl)phenoxy]-3-pyridinyl}methyl)-*N*,*N*-dimethylamine.

N,N-dimethyl-N-({3-[4-(methylsulfanyl)phenoxy]-4-pyridinyl}methyl)amine,

N-{[4-(2,3-dihydro-1-benzothien-5-yloxy)-3-pyridinyl]methyl}-N,N-dimethylamine,

N-({4-[3-Methoxy-4-(methylsulfanyl)phenoxy]-3-pyridinyl}methyl)-*N*,*N*-dimethylamine,

N,N-dimethyl-N-({4-[4-(trifluoromethyl)phenoxy]-3-pyridinyl}methyl)amine,

N,N-dimethyl-N-({4-[4-(methylsulfanyl)phenoxy]-3-pyridinyl}methyl)amine,

N,N-dimethyl-*N*-({4-[3-methyl-4-(methylsulfanyl)phenoxy]-3-pyridinyl}-methyl)amine.

- 14. (Currently amended) A composition comprising a compound of formula (I) of any one of claims 1, and 3-13, or pharmaceutically acceptable salts, solvates or polymorphs thereof, and a pharmaceutically acceptable diluent or carrier.
- 15. (Currently Amended) A therapeutic method of treating or preventing premature ejaculation comprising administering a therapeutically effective amount of a compound of formula (I) of any one of claims 1, and 3-13, or a pharmaceutically acceptable salt, solvate or polymorph thereof to a subject having a need of treatment or prevention of premature ejaculation.

16-18. (Cancelled)

- 19. (Previously presented) The compound N-methyl-N-({3-[3-methyl-4-(methylsulfanyl)phenoxy]-4-pyridinyl}-methyl)amine or a pharmaceutically acceptable salt thereof.
 - 20. (Previously presented) The tartrate salt of the compound of claim 19.

- 21. (Previously presented) The compound N-methyl-N-({3-[4-(methylsulfanyl)phenoxy]-4-pyridinyl}-methyl)amine or a pharmaceutically acceptable salt thereof.
 - 22. (Previously presented) The tartrate salt of the compound of claim 21.